

Gregory D Hawkins

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/11461457/publications.pdf>

Version: 2024-02-01

18
papers

2,954
citations

623574

14
h-index

839398

18
g-index

20
all docs

20
docs citations

20
times ranked

2286
citing authors

#	ARTICLE	IF	CITATIONS
1	Parametrized Models of Aqueous Free Energies of Solvation Based on Pairwise Descreening of Solute Atomic Charges from a Dielectric Medium. <i>The Journal of Physical Chemistry</i> , 1996, 100, 19824-19839.	2.9	828
2	Pairwise solute descreening of solute charges from a dielectric medium. <i>Chemical Physics Letters</i> , 1995, 246, 122-129.	1.2	648
3	Model for Aqueous Solvation Based on Class IV Atomic Charges and First Solvation Shell Effects. <i>The Journal of Physical Chemistry</i> , 1996, 100, 16385-16398.	2.9	358
4	Extension of the platform of applicability of the SM5.42R universal solvation model. <i>Theoretical Chemistry Accounts</i> , 1999, 103, 9-63.	0.5	197
5	Universal Quantum Mechanical Model for Solvation Free Energies Based on Gas-Phase Geometries. <i>Journal of Physical Chemistry B</i> , 1998, 102, 3257-3271.	1.2	166
6	Density functional solvation model based on CM2 atomic charges. <i>Journal of Chemical Physics</i> , 1998, 109, 9117-9133.	1.2	120
7	Improved methods for semiempirical solvation models. <i>Journal of Computational Chemistry</i> , 1995, 16, 422-440.	1.5	119
8	A universal model for the quantum mechanical calculation of free energies of solvation in non-aqueous solvents. <i>Theoretical Chemistry Accounts</i> , 1997, 98, 85-109.	0.5	99
9	Universal reaction field model based on ab initio Hartree-Fock theory. <i>Chemical Physics Letters</i> , 1998, 288, 293-298.	1.2	93
10	Parametrized Model for Aqueous Free Energies of Solvation Using Geometry-Dependent Atomic Surface Tensions with Implicit Electrostatics. <i>Journal of Physical Chemistry B</i> , 1997, 101, 7147-7157.	1.2	76
11	omnisol: a Fast Prediction of Free Energies of Solvation and Partition Coefficients. <i>Journal of Organic Chemistry</i> , 1998, 63, 4305-4313.	1.7	73
12	Universal solvation model based on conductor-like screening model. , 2000, 21, 340-366.		54
13	Prediction of Vapor Pressures from Self-Solvation Free Energies Calculated by the SM5 Series of Universal Solvation Models. <i>Journal of Physical Chemistry B</i> , 2000, 104, 4726-4734.	1.2	40
14	Solvation Modeling in Aqueous and Nonaqueous Solvents. <i>ACS Symposium Series</i> , 1994, , 24-49.	0.5	36
15	Universal Solvation Models. <i>ACS Symposium Series</i> , 1998, , 201-219.	0.5	19
16	Modeling Free Energies of Solvation and Transfer. <i>ACS Symposium Series</i> , 1998, , 285-300.	0.5	11
17	Modeling The Effect of Solvation on Structure, Reactivity, and Partitioning of Organic Solutes: Utility in Drug Design. <i>The IMA Volumes in Mathematics and Its Applications</i> , 1999, , 51-72.	0.5	10
18	New Tools for Rational Drug Design. <i>ACS Symposium Series</i> , 1999, , 121-140.	0.5	0